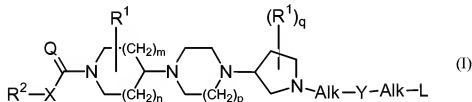


Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

1. (Previously presented) A compound according to the general Formula (I)



the pharmaceutically acceptable acid or base addition salts thereof, the stereochemically isomeric forms thereof, the *N*-oxide form thereof and prodrugs thereof, wherein :

- n is an integer, equal to 1;
- m is an integer, equal to 1;
- p is an integer equal to 1 or 2;
- q is an integer equal to 0;
- Q is O;
- X is a covalent bond;
- each R³ independently from each other, is hydrogen or alkyl ;
- each R¹ independently from each other, is selected from the group of Ar¹, Ar¹-alkyl and di(Ar¹)-alkyl ;
- R² is Ar²;
- Y is a covalent bond or a bivalent radical of formula -C(=O)-, -SO₂-, >C=CH-R or >C=N-R, wherein R is H , CN or nitro ;

- p>each Alk represents, independently from each other, a covalent bond ; a bivalent straight or branched, saturated or unsaturated hydrocarbon radical having from 1 to 6 carbon atoms ; or a cyclic saturated or unsaturated hydrocarbon radical having from 3 to 6 carbon atoms ; each radical optionally substituted on one or more carbon atoms with one or more , phenyl, halo, cyano, hydroxy, formyl and amino radicals;
- L is selected from the group of hydrogen, alkyl, alkyloxy, alkyloxyalkyloxy, alkylcarbonyloxy, alkyloxycarbonyl, mono- and di(alkyl)amino, mono- and di(alkyloxycarbonyl)amino, mono- and di(alkylcarbonyl)amino, mono- and di(Ar³)amino, mono- and di(Ar³alkyl)amino, mono- and di(Het²)amino, mono- and di(Het²alkyl)amino, alkylsulfanyl, adamantyl, Ar³, Ar³-oxy, Ar³carbonyl, Het², Het-oxy and Het²carbonyl;
- Ar¹ is phenyl, optionally substituted with 1, 2 or 3 substituents, each independently from each other, selected from the group of halo, alkyl, cyano, aminocarbonyl and alkyloxy ;
- Ar² is naphthalenyl or phenyl, each optionally substituted with 1, 2 or 3 substituents, each independently from each other, selected from the group of halo, nitro, amino, mono- and di(alkyl)amino, cyano, alkyl, hydroxy, alkyloxy, carboxyl, alkyloxycarbonyl, aminocarbonyl and mono- and di(alkyl)aminocarbonyl ;
- Ar³ is naphthalenyl or phenyl, optionally substituted with 1, 2 or 3 substituents, each independently from each other, selected from the group of alkyloxy, Ar¹carbonyloxyalkyl, Ar¹alkyloxycarbonyl, Ar¹alkyloxyalkyl, alkyl, halo, hydroxy, pyridinyl, morpholinyl, pyrrolidinyl, imidazo[1,2-*a*]pyridinyl, morpholinylcarbonyl, pyrrolidinylcarbonyl, amino and cyano;
- Het² is a monocyclic heterocyclic radical selected from the group of pyrrolidinyl, dioxolyl, imidazolidinyl, pyrazolidinyl, piperidinyl, morpholinyl, dithianyl, thiomorpholinyl, piperazinyl, imidazolidinyl, tetrahydrofuranyl, 2H-pyrrolyl, pyrrolinyl, imidazolinyl, pyrazolinyl, pyrrolyl, imidazolyl, pyrazolyl, triazolyl, furanyl, thienyl, oxazolyl, dioxazolyl, oxazolidinyl, isoxazolyl, thiazolyl, thiadiazolyl, isothiazolyl, pyridinyl, pyrimidinyl,

pyrazinyl, pyridazinyl and triazinyl ;
 or a bicyclic heterocyclic radical selected from the group of 2,3-dihydro-benzo[1,4]dioxine, octahydro-benzo[1,4]dioxine, benzopiperidinyl, quinolinyl, quinoxalinyl, indolyl, isoindolyl, chromanyl, benzimidazolyl, imidazo[1,2-*a*]pyridinyl, benzoxazolyl, benzisoxazolyl, benzothiazolyl, benzisothiazolyl, benzofuranyl or benzothienyl ;
 or the tricyclic heterocyclic radical 8,9-dihydro-4*H*-1-oxa-3,5,7a-triazacyclopenta[*f*]azulenyl ; each radical may optionally be substituted with one or more radicals selected from the group of Ar¹, Ar¹alkyl, Ar¹alkoxyalkyl, halo, hydroxy, alkyl, piperidinyl, pyrrolyl, thienyl, oxo, alkyloxy, alkylcarbonyl, Ar¹carbonyl, mono- and di(alkyl)aminoalkyl, alkyloxyalkyl and alkyloxycarbonyl; and

alkyl is a straight or branched saturated hydrocarbon radical having from 1 to 6 carbon atoms or a cyclic saturated hydrocarbon radicals having from 3 to 6 carbon atoms ; optionally substituted on one or more carbon atoms with one or more radicals selected from the group of phenyl, halo, cyano, oxo, hydroxy, formyl and amino.

2. (Previously amended) A compound according to claim 1 wherein :

R¹ is Ar¹-alkyl;

R² is Ar²;

Y is a covalent bond or a bivalent radical of formula -C(=O)-, -SO₂-, >C=CH-R or >C=N-R, wherein R is CN or nitro ;

each Alk represents, independently from each other, a covalent bond ; a bivalent straight or branched, saturated hydrocarbon radical having from 1 to 6 carbon atoms ; or a cyclic saturated hydrocarbon radical having from 3 to 6 carbon atoms ; each radical optionally substituted on one or more carbon atoms with one or more phenyl, halo and hydroxy radicals;

L is selected from the group of hydrogen, alkyl, alkyloxy, alkyloxyalkyloxy, alkylcarbonyloxy, mono- and di(alkyl)amino, mono- and di(alkyloxycarbonyl)amino, mono- and di(alkylcarbonyl)amino, mono- and di(Ar³)amino, mono- and di(Ar³alkyl)amino, mono- and di(Het²alkyl)amino,

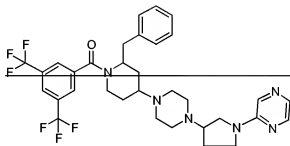
- alkylsulfanyl, adamantyl, Ar³, Het² and Het²carbonyl;
- Ar¹ is phenyl, optionally substituted with 1 or 2 halo radicals ; Ar² is naphthalenyl or phenyl, each optionally substituted with 1, 2 or 3 substituents, each independently from each other, selected from the group of halo, alkyl and alkyloxy;
- Ar³ is naphthalenyl or phenyl, optionally substituted with 1, 2 or 3 substituents, each independently from each other, selected from the group of alkyloxy, Ar¹alkyloxycarbonyl, Ar¹alkyloxyalkyl, alkyl, halo and cyano;
- Het² is a monocyclic heterocyclic radical selected from the group of pyrrolidinyl, dioxolyl, piperidinyl, morpholinyl, piperazinyl, tetrahydrofuranlyl, pyrrollyl, imidazolyl, pyrazolyl, furanyl, thienyl, dioxazolyl, oxazolidinyl, isoxazolyl, thiazolyl, thiadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl and pyridazinyl; or a bicyclic heterocyclic radical selected from the group of 2,3-dihydro-benzo[1,4]dioxine, octahydro-benzo[1,4]dioxine, quinoxalinyl, indolyl, chromanyl, benzimidazolyl, imidazo[1,2-*a*]pyridinyl, benzisoxazolyl, benzothiazolyl, benzofuranyl and benzothieryl ; or the tricyclic heterocyclic radical 8,9-dihydro-4*H*-1-oxa-3,5,7a-triazacyclopenta[f]azulenyl ; each radical may optionally be substituted with one or more radicals selected from the group of Ar¹, Ar¹alkyloxyalkyl, halo, alkyl, oxo, alkyloxy, alkylcarbonyl, Ar¹carbonyl, mono- and di(alkyl)aminoalkyl, alkyloxyalkyl and alkyloxycarbonyl ; and
- alkyl is a straight or branched saturated hydrocarbon radical having from 1 to 6 carbon atoms or a cyclic saturated hydrocarbon radicals having from 3 to 6 carbon atoms ; optionally substituted on one or more carbon atoms with one or more radicals selected from the group of phenyl, halo and hydroxy.

3. (Previously presented) A compound according to claim 1, wherein R¹ is Ar¹methyl and attached to the 2-position or R¹ is Ar¹ and attached to the 3-position.

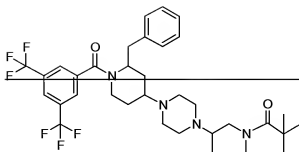
4. (Previously presented) A compound according to claim 1, wherein R²-X-C(=Q)-moiety is 3,5-di-(trifluoromethyl) phenylcarbonyl.

5. (Canceled)
6. (Previously presented) A compound according to claim 1, wherein Y is -C(=O)-.
7. (Previously presented) A compound according to claim 1, wherein Alk is a covalent bond.
8. (Previously presented) A compound according to claim 1, wherein L is Het².
9. (Currently amended) A compound according to claim 1, selected from the group consisting of:

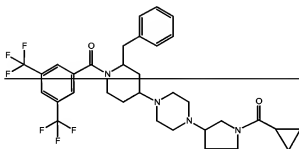
[2R-trans]-{2-benzyl-4-[4-(1-pyrazin-2-yl-pyrrolidin-3-yl)-piperazin-1-yl]-piperidin-1-yl}-(3,5-bis-trifluoromethyl-phenyl)-methanone₂



[2R-[2 α ,4 β (S)]]-1-(3-{4-[2-benzyl-1-(3,5-bis-trifluoromethyl-benzoyl)-piperidin-4-yl]-piperazin-1-yl}-pyrrolidin-1-yl)-2,2-dimethyl-propan-1-one₃

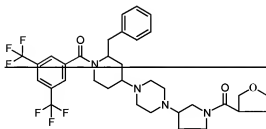


[2R-[2 α ,4 β (S*)]-{2-benzyl-4-[4-(1-cyclopropanecarbonyl-pyrrolidin-3-yl)-piperazin-1-yl]-piperidin-1-yl}-(3,5-bis-trifluoromethyl-phenyl)-methanone₄



[2R-trans]-enantiomer of {2-benzyl-4-[4-(1-cyclopropanecarbonyl-pyrrolidin-3-yl)-piperazin-1-yl]-piperidin-1-yl}-(3,5-bis-trifluoromethyl-phenyl)-methanone₄

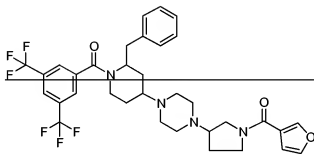
2R-trans-(2-benzyl-4-{4-[1-(tetrahydro-furan-3-carbonyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-(3,5-bis-trifluoromethyl-phenyl)-methanone₄



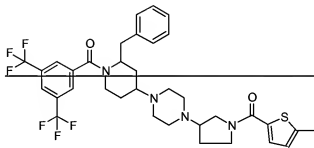
[2R-[2 α ,4 β (R(R))]]-(2-benzyl-4-{4-[1-(tetrahydro-furan-3-carbonyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-(3,5-bis-trifluoromethyl-phenyl)-methanone₄

[2R-[2 α ,4 β (S(R))]]-(2-benzyl-4-{4-[1-(tetrahydro-furan-3-carbonyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-(3,5-bis-trifluoromethyl-phenyl)-methanone₄

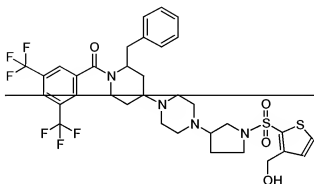
[2R-trans, R*]-(2-benzyl-4-{4-[1-(furan-3-carbonyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-(3,5-bis-trifluoromethyl-phenyl)-methanone₄



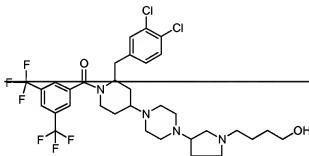
[2R-[2 α ,4 β (R))]]-(2-benzyl-4-{4-[1-(5-methyl-thiophene-2-carbonyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-(3,5-bis-trifluoromethyl-phenyl)-methanone₄



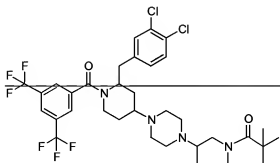
[2R-trans]-(2-benzyl-4-{4-[1-(3-hydroxymethyl-thiophene-2-sulfonyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-(3,5-bis-trifluoromethyl-phenyl)-methanone₂



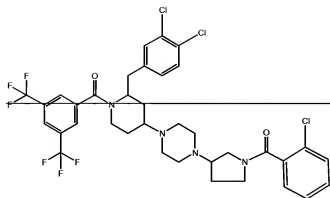
[2R-[2 α ,4 β (S)]]-(3,5-bis-trifluoromethyl-phenyl)-(2-(3,4-dichloro-benzyl)-4-{4-[1-(4-hydroxy-butyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-methanone₂



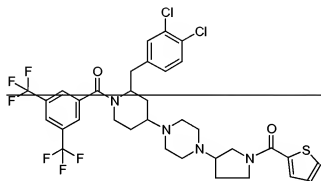
[(2R-trans),(S)]-1-(3-{4-[1-(3,5-bis-trifluoromethyl-benzoyl)-2-(3,4-dichloro-benzyl)-piperidin-4-yl]-piperazin-1-yl}-pyrrolidin-1-yl)-2,2-dimethyl-propan-1-one₂



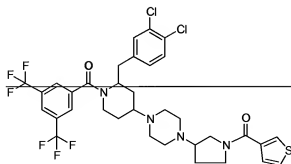
trans-(3,5-bis-trifluoromethyl-phenyl)-[4-{4-[1-(2-chloro-benzoyl)-pyrrolidin-3-yl]-piperazin-1-yl}-2-(3,4-dichloro-benzyl)-piperidin-1-yl]-methanone₂



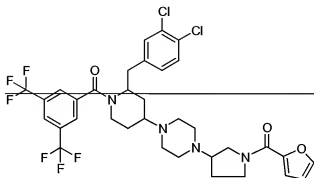
[(2R-trans),(S)]-(3,5-bis-trifluoromethyl-phenyl)-(2-(3,4-dichloro-benzyl)-4-{4-[1-(thiophene-2-carbonyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-methanone,



[(2R-trans), (R)]-(3,5-bis-trifluoromethyl-phenyl)-(2-(3,4-dichloro-benzyl)-4-{4-[1-(thiophene-3-carbonyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-methanone,

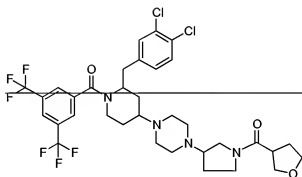


[(2R-trans), (R)]-(3,5-bis-trifluoromethyl-phenyl)-(2-(3,4-dichloro-benzyl)-4-{4-[1-(furan-2-carbonyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-methanone₄

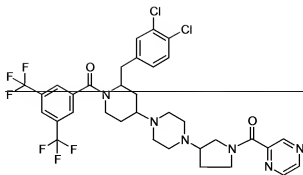


[(2R-trans), (S)]-(3,5-bis-trifluoromethyl-phenyl)-(2-(3,4-dichloro-benzyl)-4-{4-[1-(furan-2-carbonyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-methanone₄

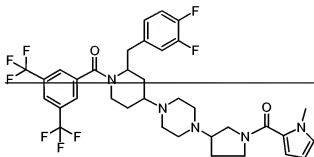
[(2R-trans), (S), (R)]-(3,5-bis-trifluoromethyl-phenyl)-(2-(3,4-dichloro-benzyl)-4-{4-[1-(tetrahydro-furan-3-carbonyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-methanone₄



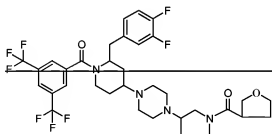
[(2R-trans), (R)]-(3,5-bis-trifluoromethyl-phenyl)-(2-(3,4-dichloro-benzyl)-4-{4-[1-(pyrazine-2-carbonyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-methanone₄



[2R-[2 α ,4 β (R*)]]-(3,5-bis-trifluoromethyl-phenyl)-(2-(3,4-difluoro-benzyl)-4-{4-[1-(1-methyl-1H-pyrrole-2-carbonyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-methanone,₂



[2R-[2 α ,4 β (R*(S*))]]-(3,5-bis-trifluoromethyl-phenyl)-(2-(3,4-difluoro-benzyl)-4-{4-[1-(tetrahydro-furan-3-carbonyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-methanone,₂

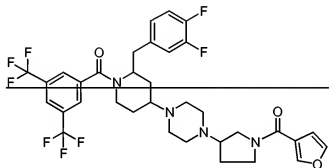


[2R-[2 α ,4 β (S*(S*))]]-(3,5-bis-trifluoromethyl-phenyl)-(2-(3,4-difluoro-benzyl)-4-{4-[1-(tetrahydro-furan-3-carbonyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-methanone₄

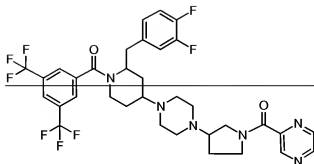
[2R-[2 α ,4 β (S*(R*))]]-(3,5-bis-trifluoromethyl-phenyl)-(2-(3,4-difluoro-benzyl)-4-{4-[1-(tetrahydro-furan-3-carbonyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-methanone₄

[2R-[2 α ,4 β (R*(R*))]]-(3,5-bis-trifluoromethyl-phenyl)-(2-(3,4-difluoro-benzyl)-4-{4-[1-(tetrahydro-furan-3-carbonyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-methanone₄

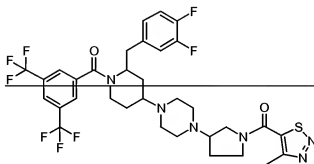
[2R-[2 α ,4 β (S*)]]-(3,5-Bis-trifluoromethyl-phenyl)-(2-(3,4-difluoro-benzyl)-4-{4-[1-(furan-3-carbonyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-methanone₄



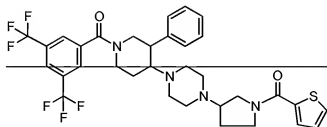
[2R-[2 α ,4 β (S*)]]-(3,5-Bis-trifluoromethyl-phenyl)-(2-(3,4-difluoro-benzyl)-4-{4-[1-(pyrazine-2-carbonyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-methanone₂



[2R-[2 α ,4 β (S*)]]-(3,5-Bis-trifluoromethyl-phenyl)-(2-(3,4-difluoro-benzyl)-4-{4-[1-(4-methyl-[1,2,3]thiadiazole-5-carbonyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-methanone, and

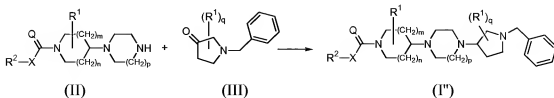


cis-(3,5-Bis-trifluoromethyl-phenyl)-(3-phenyl-4-{4-[1-(thiophene-2-carbonyl)-pyrrolidin-3-yl]-piperazin-1-yl}-piperidin-1-yl)-methanone,

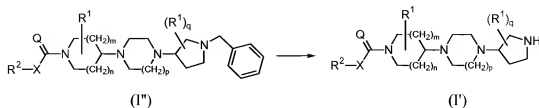


10. (Cancelled)

11. (Canceled)
12. (Canceled)
13. (Currently amended) A method for treating schizophrenia, emesis, anxiety, depression, irritable bowel syndrome (IBS), circadian rhythm disturbances, pain, neurogenic inflammation, asthma, micturition disorders ~~such as urinary incontinence~~ and nociception in a subject in need thereof comprising administering to the subject a therapeutically effective amount of a compound according to claim 1.
14. (Previously presented) A pharmaceutical composition comprising a pharmaceutically acceptable carrier and, as active ingredient, a therapeutically effective amount of a compound according to claim 1.
15. (Previously presented) A process for preparing a pharmaceutical composition as claimed in claim 14, wherein a pharmaceutically acceptable carrier is intimately mixed with a therapeutically effective amount of a compound as claimed claim 1.
16. (Original) A process for the preparation of a compound of Formula (I'') in which an intermediate compound of Formula (II) is reacted with an intermediate compound of Formula (III), wherein the radicals R^2 , X, Q, R^1 , m, n, p and q are as defined in claim 1.



17. (Withdrawn) A process for the preparation of a compound of Formula (I') in which a final compound of Formula (I'') is reductively hydrogenated, wherein the radicals R^2 , X, Q, R^1 , m, n, p and q are as defined in claim 1.



18. (Withdrawn) A process for the preparation of a compound according to Formula (I') comprising the consecutive steps of
- 1) obtaining a compound of Formula (I'') according to claim 16;
 - 2) obtaining a compound of Formula (I') according to claim 17.